

Study of the cluster splitting algorithm for the EMC

Qing Pu^{1,2}, Guang Zhao², Chunxu Yu¹, Shengsen Sun²

¹Nankai University

²Institute of High Energy Physics

PANDA Collaboration Meeting, 20/3

27/10/2020

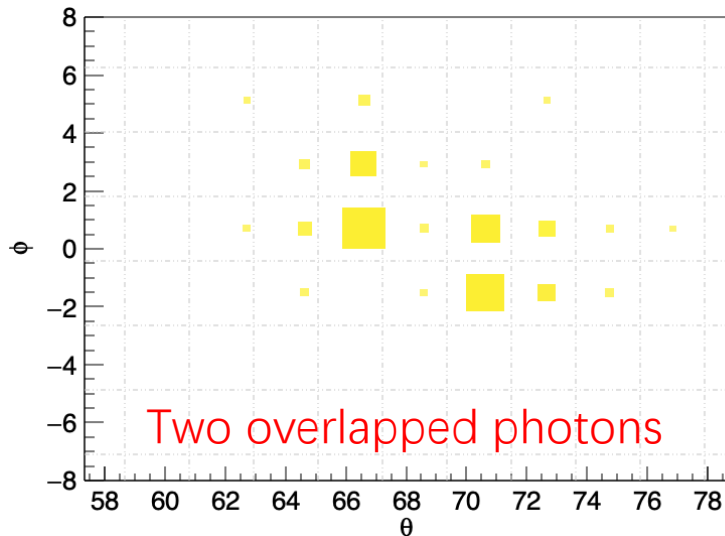
Outline

- Introduction
- Study of the cluster-splitting algorithm
- Preliminary results
- Summary

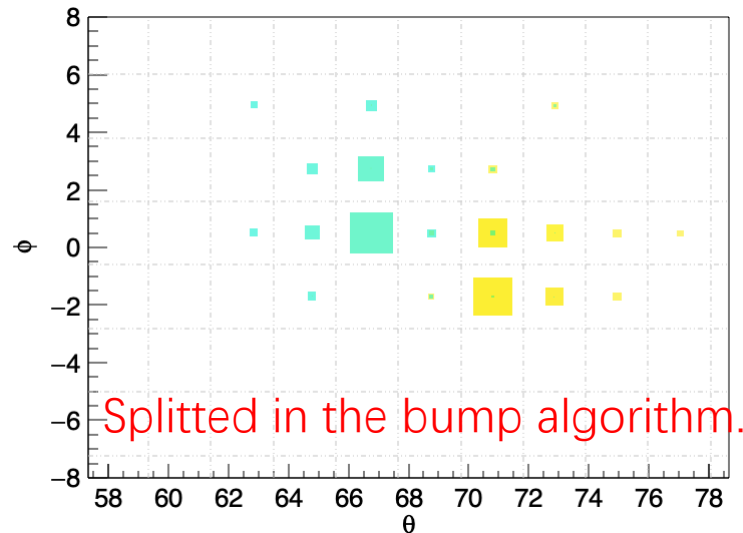
Introduction

- Cluster-splitting is an important algorithm in EMC reconstruction
- The purpose of the cluster-splitting is to separate clusters that are close to each other.
- An MC truth tool is developed to obtain the energy-fraction information in overlapped clusters (see my talk in the last collaboration meeting)
- In this work, we will study the cluster-splitting algorithm using the tool.

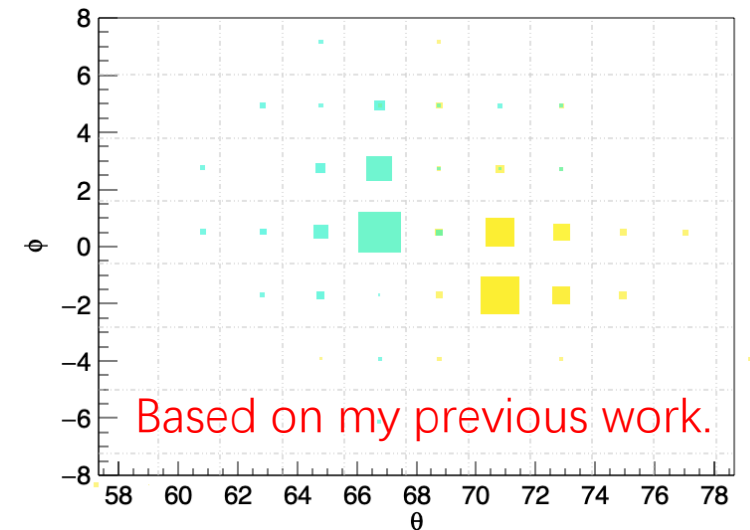
Energy distribution of Cluster



Energy distribution of Bump

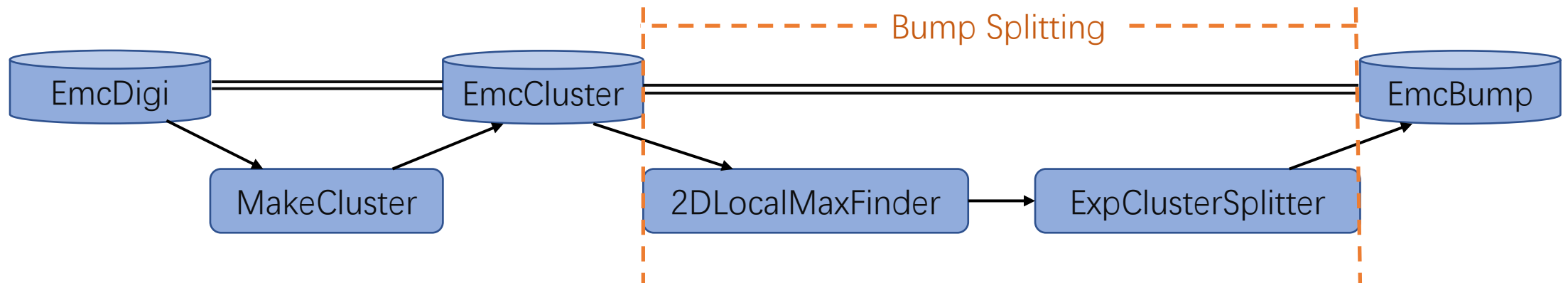


Energy distribution of MC-truth



EMC reconstruction overview

1. Cluster finding: a contiguous area of crystals with energy deposit.
2. The bump splitting
 - Find the local maximum: Preliminary split into seed crystal information
 - Update energy/position iteratively
 - The spatial position of a bump is calculated via a center-of-gravity method
 - The crystal weight for each bump is calculated by a formula.



Update energy/position iteratively

- Initialization:
Place the bump center at the seed crystal.

- Iteration:
1. Traverse all digis to calculate w_i .

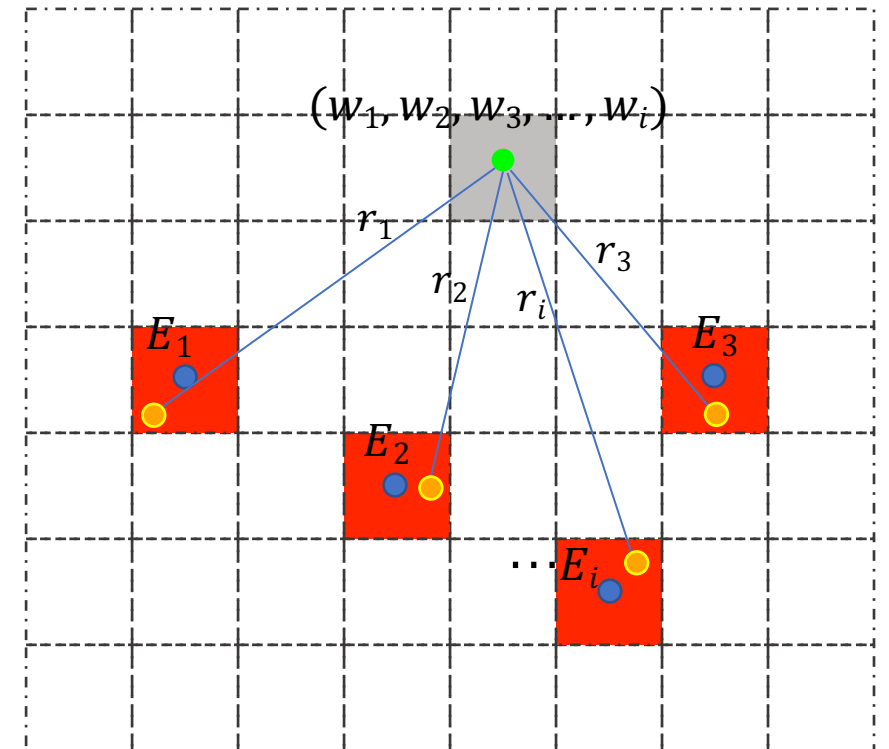
$$w_i = \frac{E_i \exp(-2.5r_i/r_m)}{\sum_j E_j \exp(-2.5r_j/r_m)}$$

i or j : different seed crystals
 r_m = Moliere radius

- 2. Update the position of the bump center.

- 3. Loop over 1 & 2 until the bump center stays stable within a tolerance of 1 mm or the number of iterations exceeds the maximum number of iterations.

- position of the crystal to be calculated
- position of seed crystal
- position of the bump center



The weighting formula

Energy deposition from a seed:

$$E_{ci}(r_{ci}) = E_i \exp(-2.5r_{ci}/r_m)$$

the two cases on the right have the same calculation result.
But the actual situation is not like this.

E_i depends not only on r_{ci} , but also on the distance from the bump center to the seed crystal r_i .

Symbol Description:

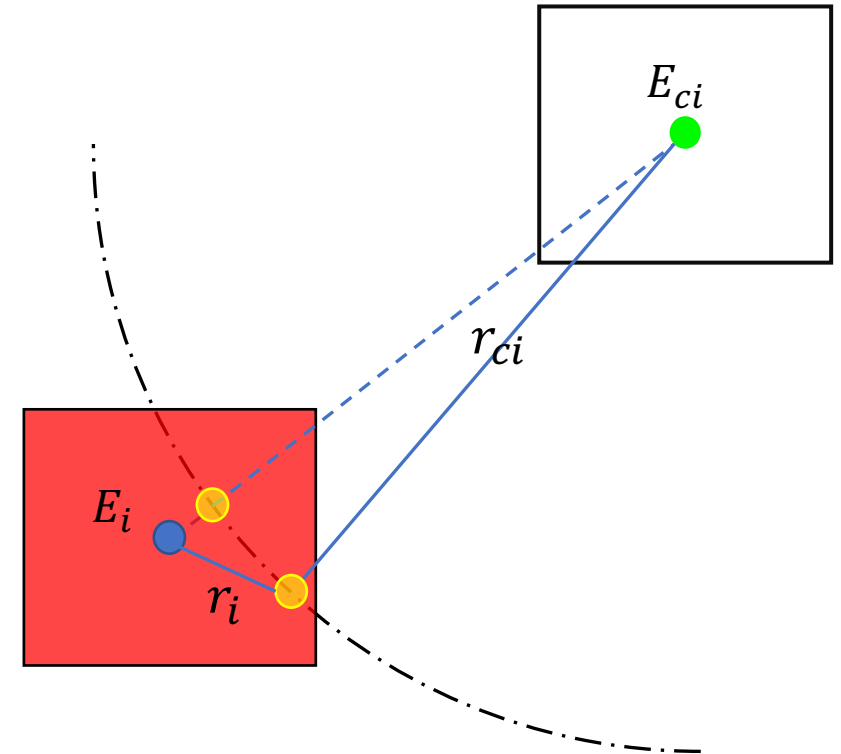
c means the current calculated crystal.

i is used to mark different seed crystals. ($i = 1, 2, 3, \dots, N_{seed}$)

E_{ci} : the deposition energy of the photon corresponding to the seed crystal i in the current crystal.

E_i : the energy of the seed crystal.

- position of the crystal to be calculated
- position of seed crystal
- position of the bump center



The weighting formula update (I)

Assuming that the bump center is located in the center of the seed crystal, there is the following function

$$E = E_s f(r)$$

E_s : the energy of the seed crystal. s = energy source

It can be obtained when the bump center is not in the center of the seed crystal :

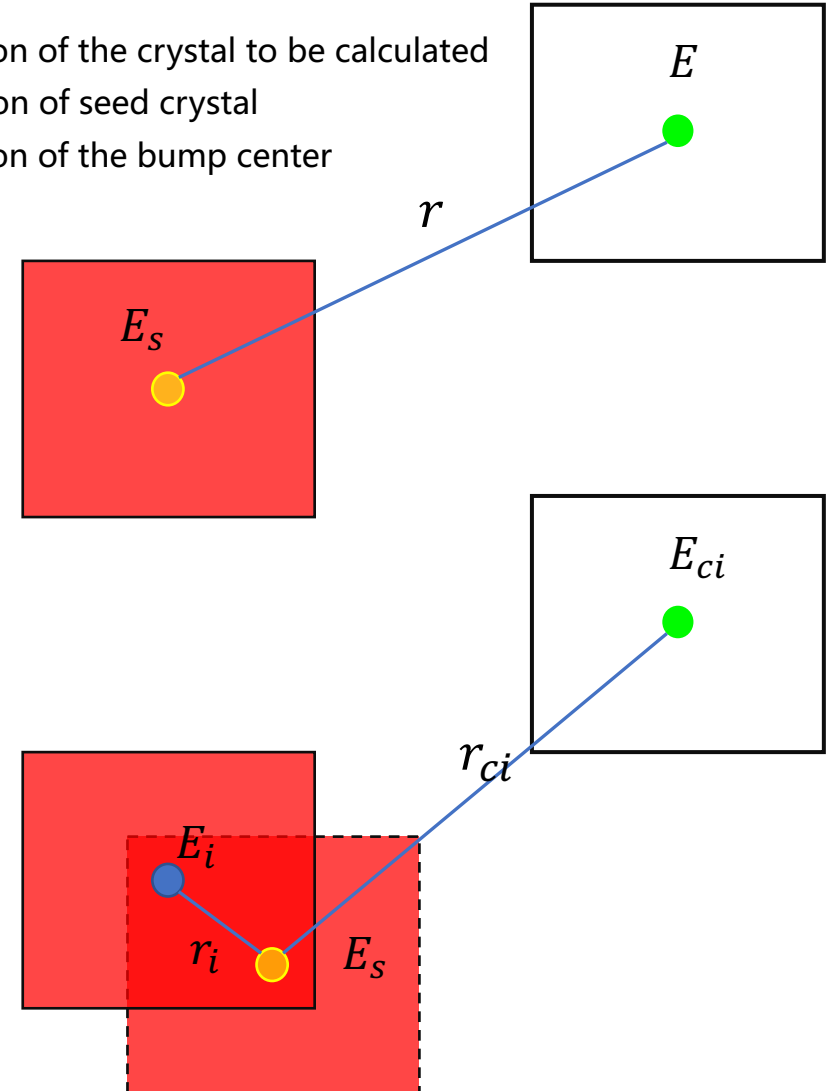
$$E_{ci} = E_s f(r_{ci}), \quad E_i = E_s f(r_i)$$

So ,

$$E_{ci} = E_i f(r_{ci}) / f(r_i)$$

- Compared with the algorithm in the library, the position information of the shower center in the seed crystal is increased.

- position of the crystal to be calculated
- position of seed crystal
- position of the bump center



The weighting formula update (II)

Assuming a certain distance, the deposition energy of crystals at different azimuths relative to the center of the shower is different.

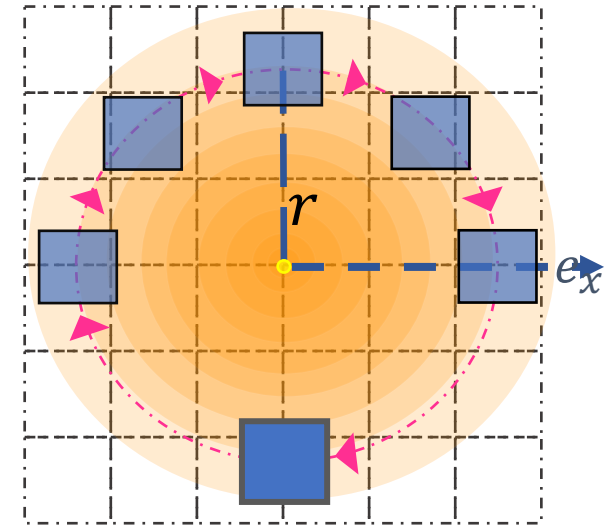
Define the azimuth as α .

The calculation formula is rewritten as

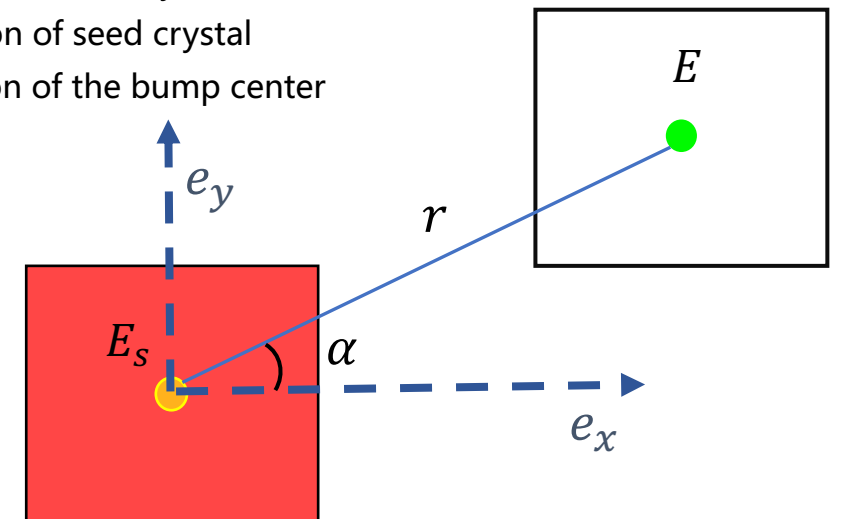
$$E = E_s f(r, \alpha)$$

So ,

$$E_{ci} = E_i f(r_{ci}, \alpha_{ci}) / f(r_i, \alpha_i)$$

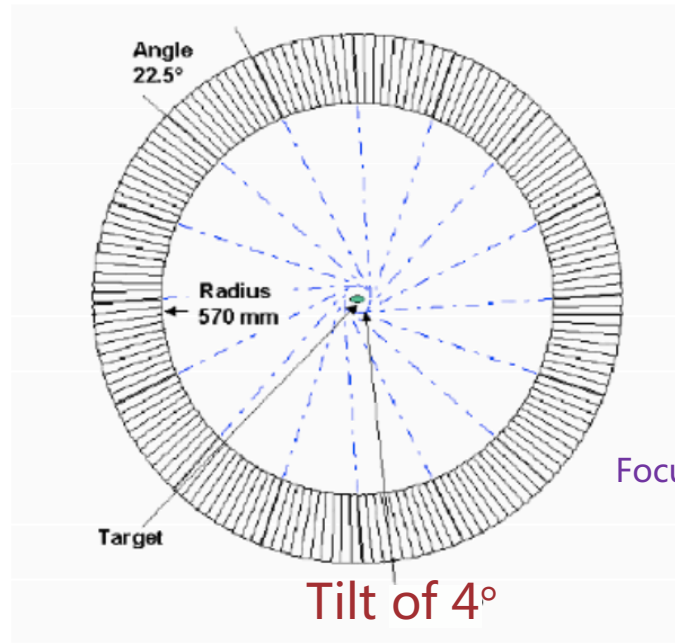
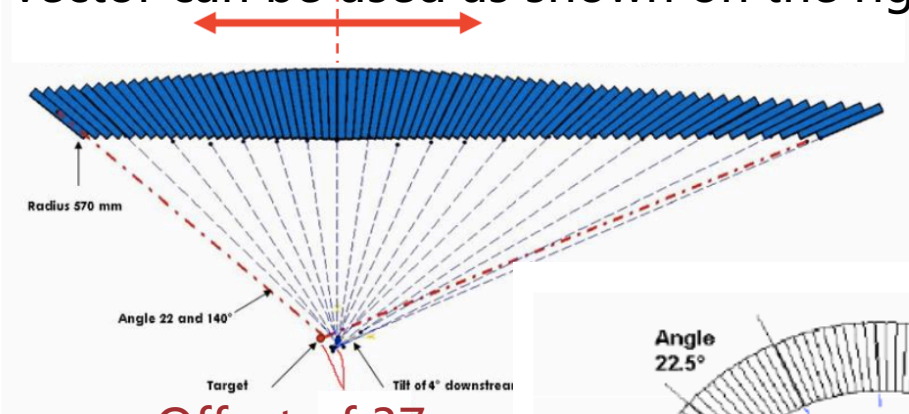


- position of the crystal to be calculated
- position of seed crystal
- position of the bump center



Calculation of angle in algorithm

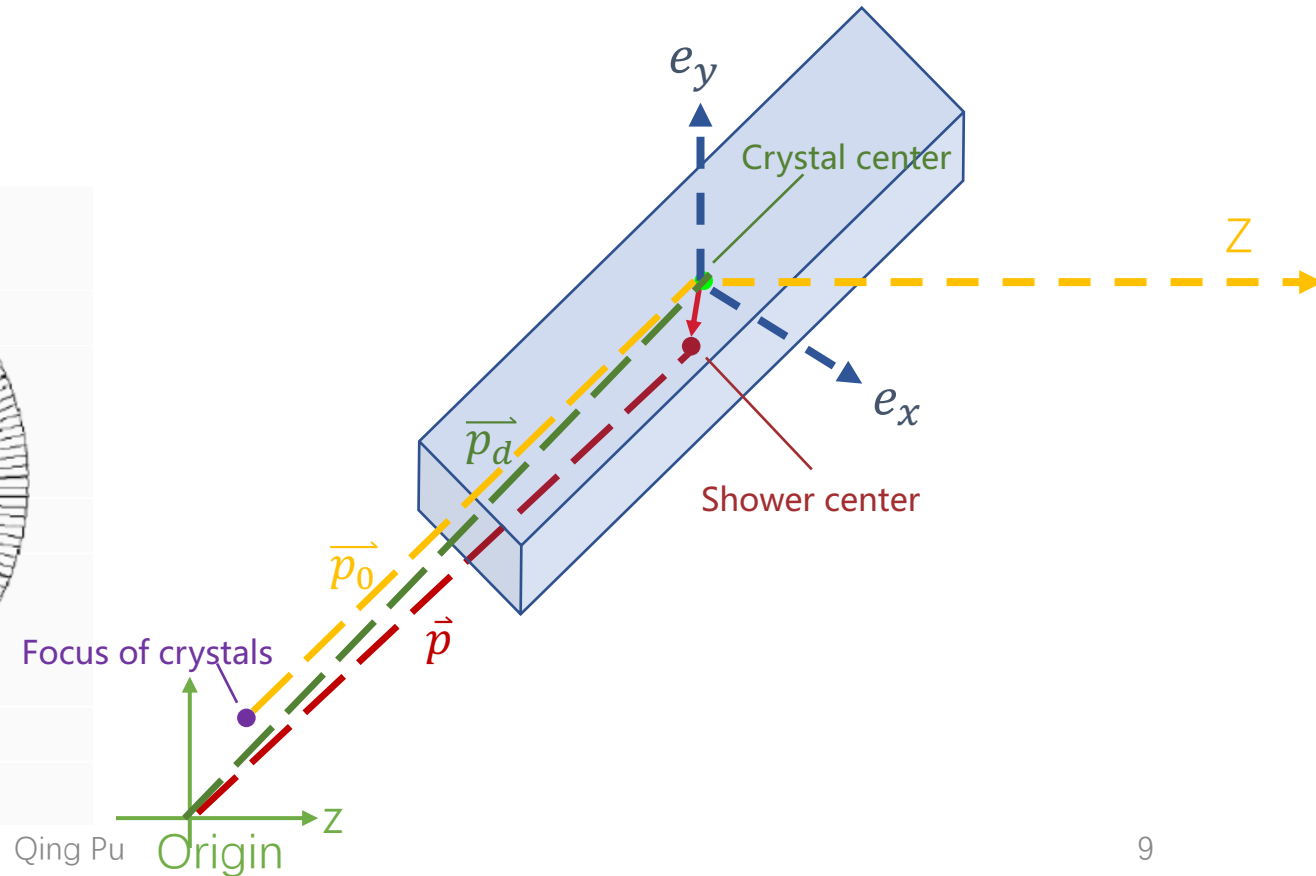
Considering the detector geometry, the normal vector can be used as shown on the right.



$$\vec{p}_0 = RotPhi(\vec{p}_d - (0, 0, 3.7), 4^\circ)$$

$$\tan\alpha = \frac{(\vec{p} - \vec{p}_0) \cdot \vec{e}_y}{(\vec{p} - \vec{p}_0) \cdot \vec{e}_x}$$

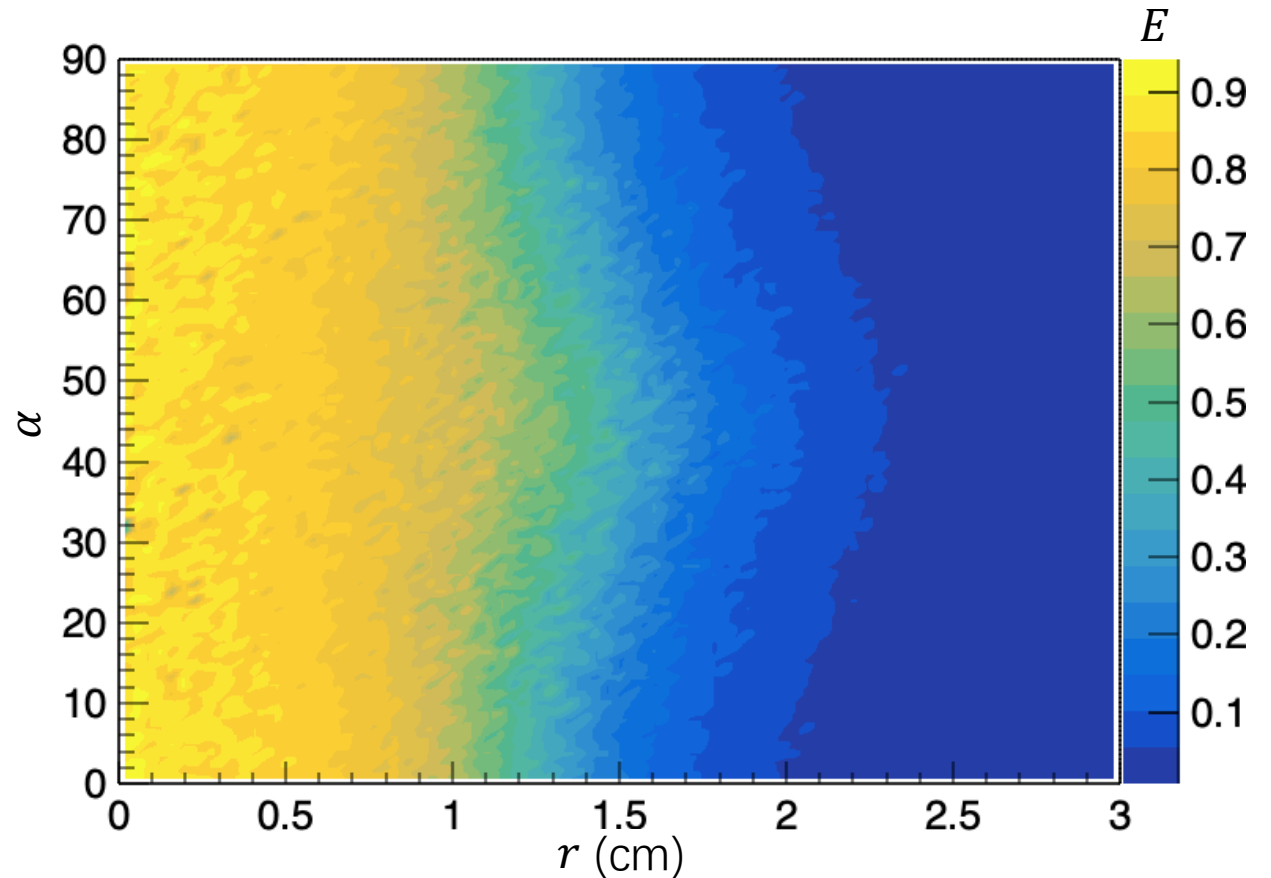
$$\left(\vec{e}_y = \frac{\vec{p}_0 \times \vec{Z}}{|\vec{p}_0|}, \vec{e}_x = \frac{\vec{p}_0 \times \vec{e}_y}{|\vec{p}_0|} \right)$$



Implement angle dependent functions

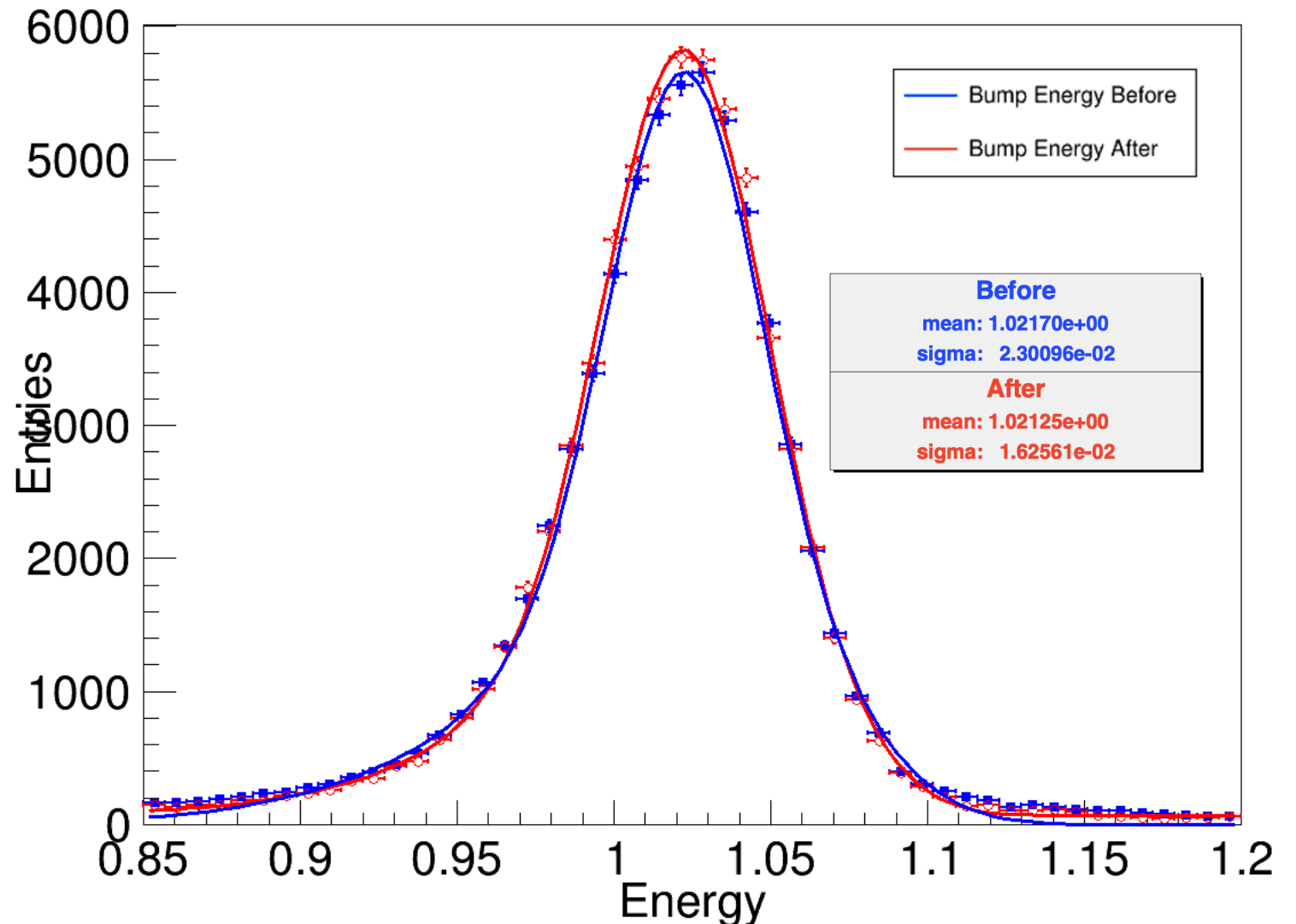
In order to get the $f(r, \alpha)$, I generated 50,000 1GeV single-photon entries and counted the distance and angle from the center of the shower.

- The energy deposition has obvious angular dependence.
- By fitting the graph on the right we can get $f(r, \alpha)$.
- Apply function $f(r, \alpha)$ to the cluster-splitting algorithm.



Preliminary Results

- Gamma Gamma (1GeV)
 - Events 50000
 - Geant3
 - Generator: Box
 - Phi(-10, 10)
 - Theta(60, 80)
-
- It can be seen that the resolution has been improved



Function used for fitting : Novosibirsk function

$$f(E) = A_S \exp(-0.5 \ln^2[1 + \Lambda \tau \cdot (E - E_0)] / \tau^2 + \tau^2)$$

Summary

- A preliminary study of the cluster-splitting algorithm is presented
- The angle and position dependency is added to the bump-weighting formula
- Test the new bump-splitting algorithm with MC-truth information. The preliminary results show slightly better resolution
- Photons of different energy and angle will be considered later.

Thank you for your attention!